

Mohamed Issam Ziane

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	A numerical optimization study of CdS and Mg _{0.125} Zn _{0.875} O buffer layers in CIGS-based solar cells using AMPS-1D package. International Journal of Modelling and Simulation, 2022, 42, 179-191.	3.3	9
2	Computational evaluation of optoelectronic, thermodynamic and electron transport properties of CuYZ ₂ (Z= S, Se and Te) chalcogenides semiconductors. Materials Chemistry and Physics, 2022, 277, 125553.	4.0	8
3	Ground-state properties of p-type delafossite transparent conducting oxides 2H-CuMO ₂ (M=Al, Sc and Tl). Journal of Applied Physics, 2019, 125, 083101.	1.9	14
4	Anisotropic optical properties of Cu ₂ ZnSn(S _x Se _{1-x}) ₄ solid solutions: First-principles calculations with TB-mBJ+U. Optik, 2021, 243, 167490.	2.9	6
5	First-Principle Computed Structural and Thermodynamic Properties of Cu ₂ ZnSn(S _x Se _{1-x}) ₄ Ternary Solid Solution. Journal of Electronic Materials, 2019, 48, 6991-7002.	2.2	5
6	First principles investigation of optoelectronic properties of ZnXP ₂ (X=Si, Ge) lattice matched with silicon for tandem solar cells applications using the mBJ exchange potential. Optik, 2018, 159, 229-244.	2.9	24
7	Optoelectronic properties of the new quaternary chalcogenides Zn ₂ CuInTe ₄ and Cd ₂ CuInTe ₄ : Ab-initio study. Optik, 2018, 157, 248-258.	2.9	6
8	The reciprocal correlation between magnetic and structural, electronic, optical properties of DMS of Zn _{1-x} SMnx. Optik, 2018, 168, 901-912.	2.9	5
9	Properties of Undoped and (Al, In) Doped ZnO Thin Films Prepared by Ultrasonic Spray Pyrolysis for Solar Cell Applications. Journal of Nano- and Electronic Physics, 2018, 10, 02036-1-02036-5.	0.5	7
10	First Principles Prediction of Structural, Electronic and Optical Properties of Zinc Blende In _x Ga _{1-x} P Alloys. Journal of Nanoelectronics and Optoelectronics, 2017, 12, 216-223.	0.5	1
11	Full-potential calculations of structural and optoelectronic properties of cubic indium gallium arsenide semiconductor alloys. Optik, 2016, 127, 9280-9294.	2.9	27
12	First-principles investigation of the optical properties for rocksalt mixed metal oxide Mg _{1-x} Zn _x O. Materials Chemistry and Physics, 2016, 182, 182-189.	4.0	28
13	First Principal Calculations of Optical Properties of InGa ₂ Using in Solar Cells Applications. Applied Condition Monitoring, 2015, , 179-187.	0.4	1
14	First principles study of structural, electronic and optical properties of indium gallium nitride arsenide lattice matched to gallium arsenide. Materials Science in Semiconductor Processing, 2015, 30, 181-196.	4.0	31
15	First-principles prediction of the structural and electronic properties of zinc blende Ga _x As _{1-x} alloys. Materials Science in Semiconductor Processing, 2013, 16, 1138-1147.	4.0	17